## WEST

# Freeform Search

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### **Search History**

<b>DB Name</b>	Query	Hit Count S	Set Name
ALL	153 same 147	18	<u>L54</u>
ALL	140 near10 152	1212	<u>L53</u>
ALL	phospho\$5	377624	<u>L52</u>
ALL	149 and 12	4	<u>L51</u>
ALL	148 and 12	130	<u>L50</u>
ALL	148 same phospho\$5	227	<u>L49</u>
ALL	-147 same 140	3384	<u>L48</u>
ALL	conduct\$5	1746061	<u>L47</u>
ALL	140 near 10 141	71	<u>L46</u>
ALL	complex near10 140 near10 141	3	<u>L45</u>
ALL	142 and 12	90	<u>L44</u>
ALL	142 same 139	0	<u>L43</u>
ALL	140 same 141	732	<u>L42</u>
ALL	\$8phosphonic	27476	<u>L41</u>
ALL	dianiline or aniline or dithiodianiline or dimethylaniline or diethylaniline or bisaniline	67033	<u>L40</u>

ALL	conductivity	269810	<u>L39</u>
ALL	137 and 12	10	<u>L38</u>
ALL	(stillbene or thiopene or carbazole or pyrene or pyrrole or phenanthrene) near5 carboxaldehyde	264	<u>L37</u>
ALL	135 and 12	56	<u>L36</u>
ALL	(methoxy or hydroxy) near5 naphthaldehyde	434	<u>L35</u>
ALL	133 and 12	0	<u>L34</u>
ALL	bromo adj3 furaldehyde	29	<u>L33</u>
ALL	131 and 12	6	<u>L32</u>
ALL	(nitro or diethylamino or methoxy) near5 cinnamaldehyde	73	<u>L31</u>
ALL	(nitro or diethylamino or methoxy) near5 cinnaldehyde	0	<u>L30</u>
ALL	127 and 12	4	<u>L29</u>
ALL	l27 and l17	4	<u>L28</u>
ALL	hydroxy near5 nitrobenzaldehyde	209	<u>L27</u>
ALL	(dibenzyloxy or dimethoxy or dibromo or trimethoxy) near5 benzaldehye	0	<u>L26</u>
ALL	124 and 12	137	<u>L25</u>
ALL	(hydroxy or benzyloxy or carboxy or nitro or dihydroxy or methoxy or ethoxy or methyl) near5 benzaldehyde	4227	<u>L24</u>
ALL	122 and ink.ti.	12	<u>L23</u>
ALL	119 and 12	285	<u>L22</u>
ALL	120 and 12	2	<u>L21</u>
ALL	119 and 16	2	<u>L20</u>
ALL	cinnamaldehyde or naphthaldehyde or anthraldehyde or carboxaldehyde or dicarboxaldehyde or pyrrolecarboxaldehyde or fluorenecarboxaldehyde	8338	<u>L19</u>
ALL	117 and 12.ti.	54	<u>L18</u>
ALL	115 and 12	1007	<u>L17</u>
ALL	115 and 16	16	<u>L16</u>
ALL	benzaldehyde or nitrobenzaldehyde or trichlorobenzaldehyde or dimethoxybenzaldehyde	22129	<u>L15</u>
ALL	16 and 15	91	<u>L14</u>
ALL	(phenyl adj glycidyl adj ether) near5 15	5	<u>L13</u>
ALL	(cresyl adj glycidyl adj ether) near5 15	0	<u>L12</u>
ALL	110 and 12	129	<u>L11</u>
ALL	15 same 14	1124	<u>L10</u>
ALL	16 and 15 and 14	6	<u>L9</u>
ALL	16 and (15 same 14)	2	<u>L8</u>
ALL	16 and 13	83	<u>L7</u>

ALL	11 near5 l2	2945	<u>L6</u>
ALL	formaldehyde	119017	<u>L5</u>
ALL	"glycidyl ether" or toluenesulfonamide	18770	<u>L4</u>
ALL .	aldehyde	106811	<u>L3</u>
ALL	ink	263562	<u>L2</u>
ALL	"hot melt" or "phase change"	61701	<u>L1</u>

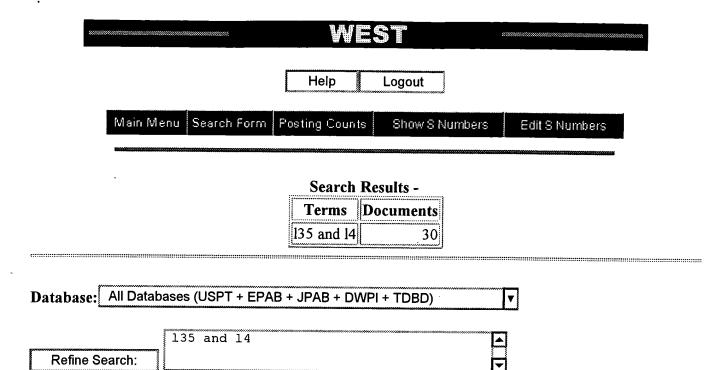
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	BRS	BRS	BRS	BRS	BRS	BRS	BRS	IS&R	Type
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	21851	235	0	1736619	549265	376727	2634	14	Hits
F	phenylphosphonic or phosphonic	dithiodianiline or dimethyldianiline or diethyldianiline or bisaniline	5 and 6	conduct\$6	complex	phospho\$5	dianiline	("528628") or ("6022910") or ("5041161") or ("6017385") or ("5041161") or ("5017385") or ("5006170") or ("5931995") or ("5939468") or ("5825391") or ("4020040") or ("5902390") or ("5876492") or ("5931995") or ("5270730") or ("6017385") or ("5270730") or ("6017385") or ("5713989")).PN.	Search Text
Derwent	R;		USPAT; USOCR; EPO; JPO; Derwent	USPAT; USOCR; EPO; JPO; Derwent	USPAT; USOCR; EPO; JPO; Derwent	USPAT; USOCR; EPO; JPO; Derwent	USPAT; USOCR; EPO; JPO; Derwent	USPAT	DBs
	2000/05/01	2000/05/01 11:28	2000/05/01 11:28	2000/05/01 11:28	2000/05/01 11:26	2000/05/01 11:26	2000/05/01 11:25	2000/05/01	Time Stamp
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Type	BRS	10 BRS I	11 BRS I	12 BRS I	13 BRS I	14 BRS I	15 BRS I	16 BRS I	
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DBs	USPAT; USOCR; EPO; JPO; Derwent	USPAT; USOCR; EPO; JPO; Derwent	USPAT; USOCR; EPO; JPO; Derwent	7	USPAT; USOCR; EPO; JPO; Derwent	USPAT; USOCR; EPO; JPO; Derwent	rt	Ct .	USPAT; USOCR; EPO;
Time Stamp	2000/05/01 11:28	2000/05/01 11:29	2000/05/01 11:29	2000/05/01 11:30	2000/05/01 11:30	2000/05/01 11:30	2000/05/01 11:31	2000/05/01 11:31	2000/05/01
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0			2000/05/01 11:49		formaldehyde	118928 f	L32 1	BRS	26 I
0			2000/05/01 11:49	4	aldehyde	106793 a	L31 1	BRS	25
0			2000/05/01 11:49	USPAT; USOCR; EPO; JPO; Derwent	ink	262070 iı	L30 2	BRS	24
0			2000/05/01 11:49		"hot melt" or phase adj2 change	65620 ".	L29 e	BRS	23
0			2000/05/01 11:46	USPAT; USOCR; EPO; JPO; Derwent	5 same 3 same 5	60 25	L28 (	BRS	22
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0			2000/05/01 11:35	USPAT; USOCR; EPO; JPO; Derwent	polyaniline '	3675 p	Ĺ25	BRS	20
0			2000/05/01 11:34	rt	aniline	53982 a	L23	BRS	19
0			2000/05/01 11:34	USPAT; USOCR; EPO; JPO; Derwent	same 3	53	L22	BRS	18
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	36 and 30			35 and 30				3/ alia 38	) }				37 and 35				29 near5 30			24 HEALD 32	7 0 0 7 7			33 nearb 32	ì			toluenesulfonamide			gryciayi adj etner	) L I.	••
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PC:11				11:51				11:51	2000/05/01			11:50	2000/05/01			11:50	2000/05/01			11:50	2000/05/01			11:56	2000/05/01			2000/05/01				2000/05/01	••
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#### Search History

<b>DB Name</b>	Query	Hit Count	Set Name
ALL	135 and 14	30	<u>L36</u>
ALL	"phenyl glycidyl ether" same formaldehyde	117	<u>L35</u>
ALL	"phenyl glycidyl ether" adj3 formaldehyde	1	<u>L34</u>
ALL	132 and 14	10	<u>L33</u>
ALL	(polyglycidyl same 118) same 12	62	<u>L32</u>
ALL	(127 same 12) and 14	12	<u>L31</u>
ALL	127 and 14	103	<u>L30</u>
ALL	128 and 14	1	<u>L29</u>
ALL	127 and 16	1	<u>L28</u>
ALL	(glycidyl adj ether) same 118	990	<u>L27</u>
ALL	(125 same 120) and 16	14	<u>L26</u>
ALL	aldehyde	106811	<u>L25</u>
ALL	123 and 16	13	<u>L24</u>
ALL	phenol adj formaldehyde	22353	<u>L23</u>
ALL	18 and 115	57	<u>L22</u>
ALL	(120 same 118) and 16	0	<u>L21</u>
ALL	vehicle	1012955	<u>L20</u>
ALL	118 and 16	91	<u>L19</u>
ALL	formaldehyde	119017	L18

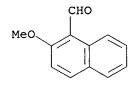
ALL	formaldehyde	119017	<u>L18</u>
ALL	114 and 115	64	<u>L17</u>
ALL	14 and 15	6109	<u>L16</u>
ALL	ink.ti.	97431	<u>L15</u>
ALL	l6 and 11	143	<u>L14</u>
ALL	16 amd 11	0	<u>L13</u>
ALL	111 and 18	15	<u>L12</u>
ALL	19 near5 110	9728	<u>L11</u>
ALL	modifie\$2 or modify	814824	<u>L10</u>
ALL	viscosity	380156	<u>L9</u>
ALL	14 and 13	486	<u>L8</u>
ALL	16 and 13	6	<u>L7</u>
ALL	14 near5 15	2945	<u>L6</u>
ALL	"hot melt" or "phase change"	61701	<u>L5</u>
ALL	ink	263562	<u>L4</u>
ALL	11 near5 12	4099	<u>L3</u>
ALL	copolymer\$1	454934	<u>L2</u>
ALL	aldehyde or formaldehyde	198612	<u>L1</u>

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ANSWER 1 OF 1 REGISTRY COPYRIGHT 2000 ACS
    104-55-2 REGISTRY
    2-Propenal, 3-phenyl- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
    Cinnamaldehyde (8CI)
OTHER NAMES:
     .beta.-Phenylacrolein
CN
     3-Phenyl-2-propenal
CN
     3-Phenyl-2-propenaldehyde
CN
     3-Phenylacrolein
CN
     3-Phenylacrylaldehyde
CN
     3-Phenylpropenal
CN
    Abion CA
CN
    Benzylideneacetaldehyde
CN
     Cassia aldehyde
CN
     Cinnamal
CN
     Cinnamic aldehyde
CN
     Cinnamyl aldehyde
CN
     Phenylacrolein
CN
     Zimtaldehyde
CN
     3D CONCORD
FS
MF
     C9 H8 O
     COM
CI
     STN Files: AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
LC
       BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS,
       CHEMINFORMRX, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DETHERM*, DRUGU,
       EMBASE, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*,
       MSDS-OHS, NAPRALERT, NIOSHTIC, PDLCOM*, PIRA, PROMT, RTECS*, SPECINFO,
       TOXLINE, TOXLIT, TULSA, ULIDAT, USAN, USPATFULL, VETU, VTB
         (*File contains numerically searchable property data)
     Other Sources: DSL**, EINECS**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
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5169 REFERENCES IN FILE CA (1967 TO DATE)
70 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
5179 REFERENCES IN FILE CAPLUS (1967 TO DATE)
5 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

ANSWER 1 OF 1 REGISTRY COPYRIGHT 2000 ACS L2 5392-12-1 REGISTRY RN 1-Naphthalenecarboxaldehyde, 2-methoxy- (9CI) (CA INDEX NAME) CN OTHER CA INDEX NAMES: 1-Naphthaldehyde, 2-methoxy- (6CI, 7CI, 8CI) CN OTHER NAMES: CN 2-Methoxy-1-naphthaldehyde 2-Methoxy-1-naphthalenecarboxaldehyde CN2-Methoxy-1-naphthylaldehyde CN 3D CONCORD FS C12 H10 O2 MF BEILSTEIN\*, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, LCSTN Files: CHEMINFORMRX, CHEMLIST, CSCHEM, HODOC\*, IFICDB, IFIPAT, IFIUDB, MEDLINE, SPECINFO, TOXLIT, USPATFULL (\*File contains numerically searchable property data) Other Sources: EINECS\*\* (\*\*Enter CHEMLIST File for up-to-date regulatory information)



142 REFERENCES IN FILE CA (1967 TO DATE)
142 REFERENCES IN FILE CAPLUS (1967 TO DATE)
4 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

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L6
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2000 ACS
RN
     1899-24-7 REGISTRY
CN
     2-Furancarboxaldehyde, 5-bromo- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
     2-Furaldehyde, 5-bromo- (6CI, 7CI, 8CI)
OTHER NAMES:
     2-Bromo-5-furancarboxaldehyde
     2-Bromofuran-5-carboxaldehyde
CN
CN
     5-Bromo-2-formylfuran
CN
     5-Bromo-2-furaldehyde
     5-Bromo-2-furancarboxaldehyde
CN
     5-Bromo-2-furfural
CN
     5-Bromofurfural
CN
FS
     3D CONCORD
MF
     C5 H3 Br O2
CI
     COM
                  BEILSTEIN*, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS,
LC
       CHEMINFORMRX, CSCHEM, HODOC*, IFICDB, IFIPAT, IFIUDB, SPECINFO, TOXLIT,
       USPATFULL
         (*File contains numerically searchable property data)
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228 REFERENCES IN FILE CA (1967 TO DATE)
229 REFERENCES IN FILE CAPLUS (1967 TO DATE)
14 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

ANSWER 1 OF 1 REGISTRY COPYRIGHT 2000 ACS L9 4521-33-9 REGISTRY RN 2-Thiophenecarboxaldehyde, 5-nitro- (6CI, 7CI, 8CI, 9CI) (CA INDEX NAME) CNOTHER NAMES:  $\hbox{$2$-Formyl-$5$-nitrothiophene}$ CN 2-Nitrothiophene-5-carboxaldehyde CN 5-Nitro-2-formylthiophene CN 5-Nitro-2-thienylaldehyde CN 5-Nitro-2-thienylcarboxaldehyde CN CN 5-Nitro-2-thiophenealdehyde 5-Nitro-2-thiophenecarboxaldehyde CN CN NTA 3D CONCORD FS C5 H3 N O3 S MF CI COM BEILSTEIN\*, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, LCSTN Files: CHEMINFORMRX, CHEMLIST, CSCHEM, IFICDB, IFIPAT, IFIUDB, TOXLIT, USPATFULL (\*File contains numerically searchable property data) Other Sources: EINECS\*\* (\*\*Enter CHEMLIST File for up-to-date regulatory information)

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190 REFERENCES IN FILE CA (1967 TO DATE)
4 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
190 REFERENCES IN FILE CAPLUS (1967 TO DATE)